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# **A Complex Wavelet representation of Error Covariances in ALADIN 3d-Var**

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- Introduction
- A complex wavelet approach in 1+1 dimensions
- Full 3d implementation: first experiments
- Conclusions

# Variational data assimilation

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- In 3 dimensions:

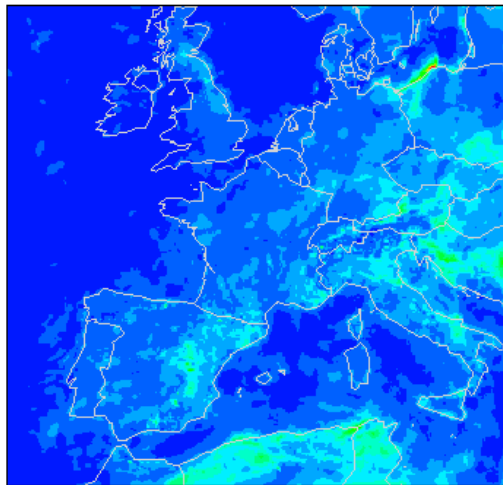
$$J(\mathbf{x}) = \frac{1}{2}(\mathbf{x} - \mathbf{x}^b)^T \mathbf{B}^{-1}(\mathbf{x} - \mathbf{x}^b) + \frac{1}{2}(\mathbf{H}\mathbf{x} - \mathbf{y})^T \mathbf{R}^{-1}(\mathbf{H}\mathbf{x} - \mathbf{y}),$$

- $\mathbf{H}$  is the **observation operator**.  $\mathbf{B}$  and  $\mathbf{R}$  are the **covariance matrices** of the background and observation errors, respectively.
- Formal exact solution:
$$\mathbf{x}^a = \mathbf{x}^b + \mathbf{B}\mathbf{H}^T (\mathbf{H}\mathbf{B}\mathbf{H}^T + \mathbf{R})^{-1} (\mathbf{y} - \mathbf{H}\mathbf{x}^b)$$
- $\mathbf{B}$  describes how one observation influences the analysis in the neighbourhood.

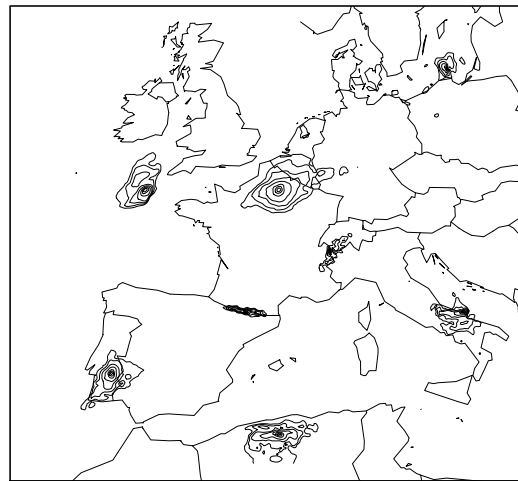
# Background Error Covariance

- The matrix  $B$  of background error covariances is vital to most assimilation methods.

T-41 variance



T-41 correlations



- Heterogeneous, anisotropic structure functions (and often noisy).
- May be estimated in many ways, either statically or day by day (ensemble).

# Simplifying $\mathbf{B}$

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- $\mathbf{B}$  is much too large as a full matrix.
- A **diagonal** matrix requires much less memory (and inverting it is trivial).
- If we represent  $\mathbf{B}$  in grid point space, the diagonal represents the error **variance** at every grid point. The correlation of errors in different points is completely ignored.

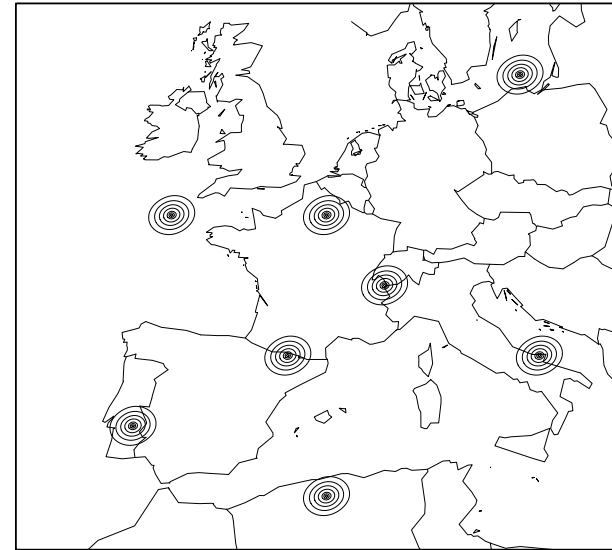
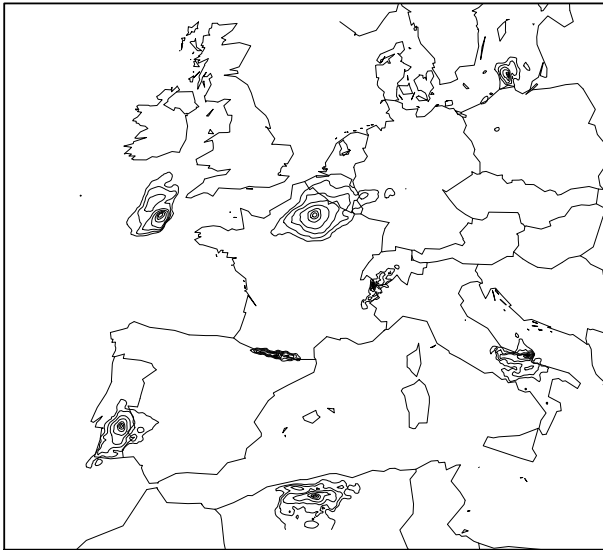
# Diagonalising B

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- We can represent our background field in **spectral** co-ordinates:  $\mathbf{x}_f = \mathbf{F}\mathbf{x}$ .
- The B matrix in Fourier space is  $\mathbf{B}_f = \mathbf{F}\mathbf{B}\mathbf{F}^*$ .
- If we diagonalise  $\mathbf{B}_f$ , the covariance matrix in grid point space becomes  $\tilde{\mathbf{B}} = \mathbf{F}^*\mathbf{B}_f\mathbf{F}$
- A diagonal matrix in Fourier space corresponds to **homogeneous** structure functions in grid space.
- One may combine these 2 representations  $\mathbf{B} = \mathbf{D}_g\mathbf{F}^*\mathbf{C}_f\mathbf{F}\mathbf{D}_g$ , where  $\mathbf{D}_g$  represents the standard deviation and  $\mathbf{C}_f$  the correlations.

# Spectral diagonalisation

- Spectral diagonalisation of  $B$  gives homogeneous correlation functions:



- One “average” structure function for the whole domain.
- Can we do better than this?

# Introducing wavelets

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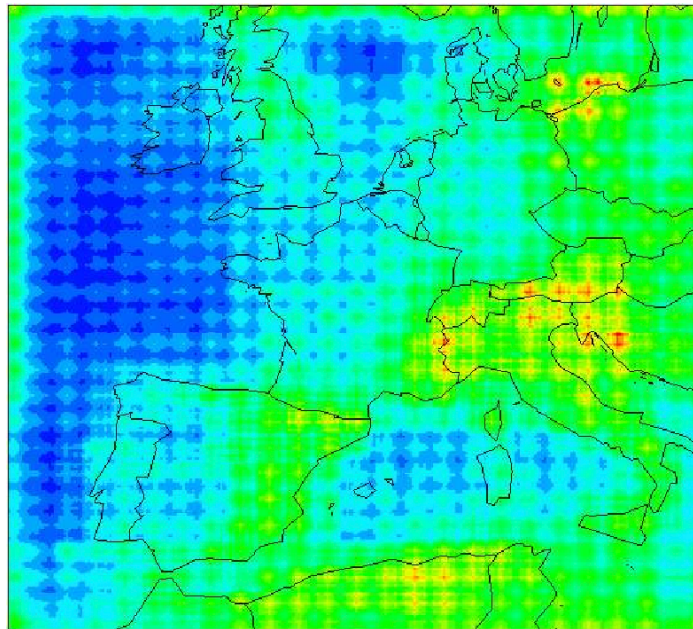
- M. Fisher (2001) introduced the idea of representing  $B$  in wavelet co-ordinates.
- For the *global* model IFS used at ECMWF he introduced a set of non-orthogonal, band-limited wavelets.
- No longer an orthogonal basis, but a **tight frame**.



# A first (naive) approach

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- Using 1 single orthonormal basis of (Meyer) wavelets, the variance (diagonal of  $\mathbf{B}$ ) in grid point space becomes:



- The wavelet transform & diagonalisation introduce unacceptable artifacts.

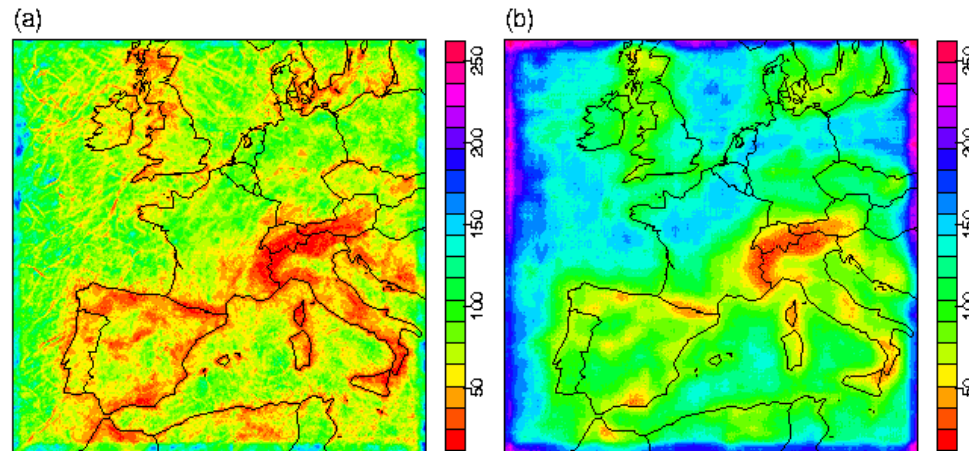
# A hybrid Meyer approach

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- In Deckmyn & Berre (2005) wavelets were combined with grid point and Fourier.
- Use every basis for its strongest points:
  - **Grid space**: strictly local (variance)
  - **Fourier space**: average correlation function
  - **Wavelet space**: local differences from average
- $\mathbf{B} = \mathbf{D}_g^* \mathbf{F}^* \mathbf{D}_f^* (\mathbf{F}^{-1})^* \mathbf{W}^* \mathbf{B}_w \mathbf{W} \mathbf{F}^{-1} \mathbf{D}_f \mathbf{F} \mathbf{D}_g,$

# A hybrid Meyer approach

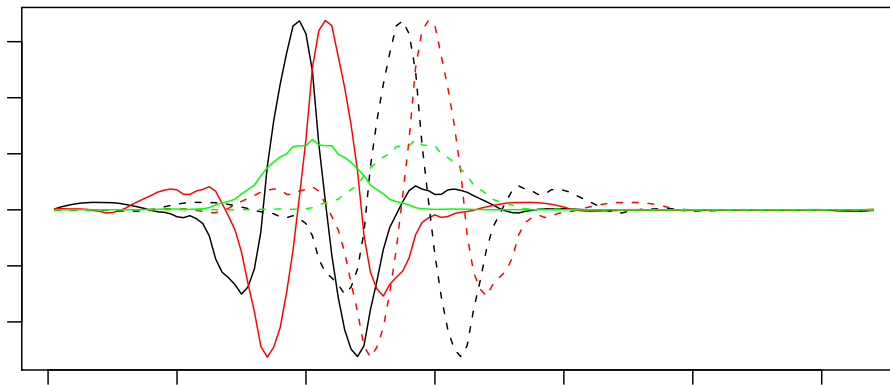
- The length scale (width of the local correlation function  $L^2 \approx \frac{-2\rho}{\nabla^2 \rho}$ ) becomes:



- In fact a clearer image than the noisy original!
- Quite cumbersome. A lot of calculations for relatively small gain.

# Complex wavelets

- Introduced by N. Kingsbury (2001).
- 2 separate (“Q-shift dual tree”) orthonormal wavelet transforms, that can be interpreted as real and imaginary components ( $\approx$  windowed  $\cos$  &  $\sin$ )



- Approximate **shift invariant**  $\rightarrow$  much less artifacts!
- Limited redundancy of 2:1.

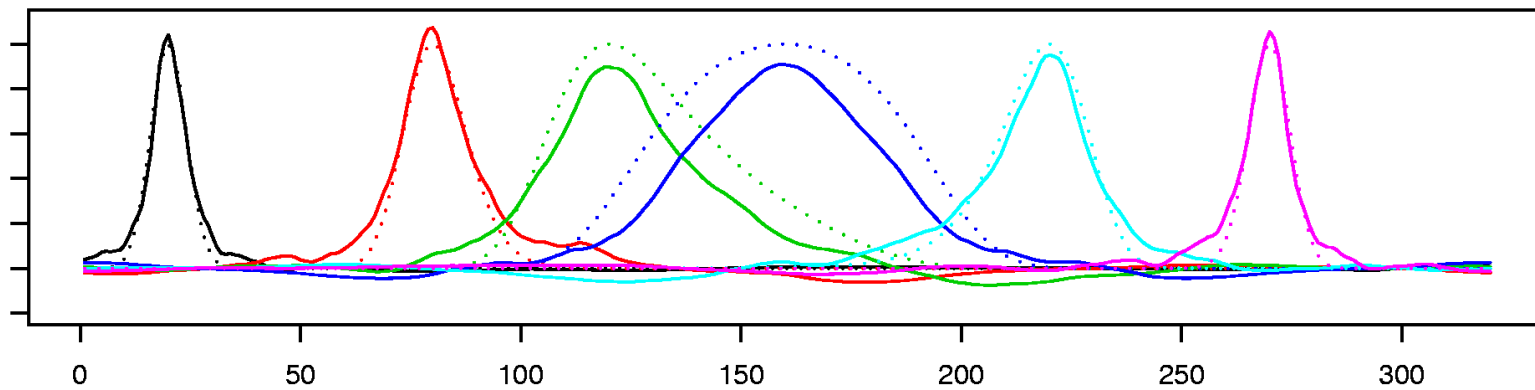
# Complex wavelets

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- The two wavelet filters are about  $1/2$  sample apart.
- To achieve this, the first stage of  $\mathbf{W}_2$  in fact uses a different filter.
- The complex wavelet transform  $\mathbf{W} = \mathbf{W}_1 + i\mathbf{W}_2$  (or, equivalently, the combination of  $\mathbf{W}_1$  and  $\mathbf{W}_2$  as a set of  $2N$  real functions) is not orthonormal, but a tight frame of multiplicity 2.
- The two wavelets are in fact the reverse of each other!
- $\mathbf{B}_w$  is a complex, hermitian  $N \times N$  matrix.

# 1D Complex wavelets

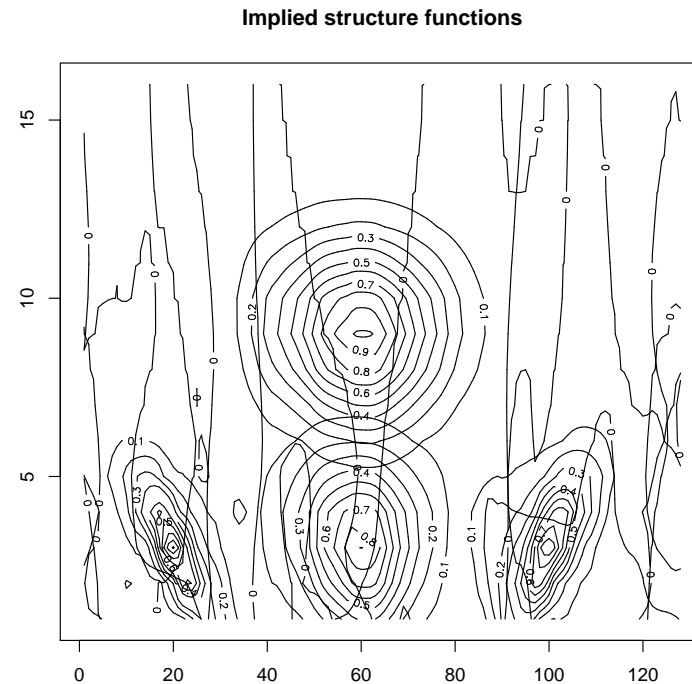
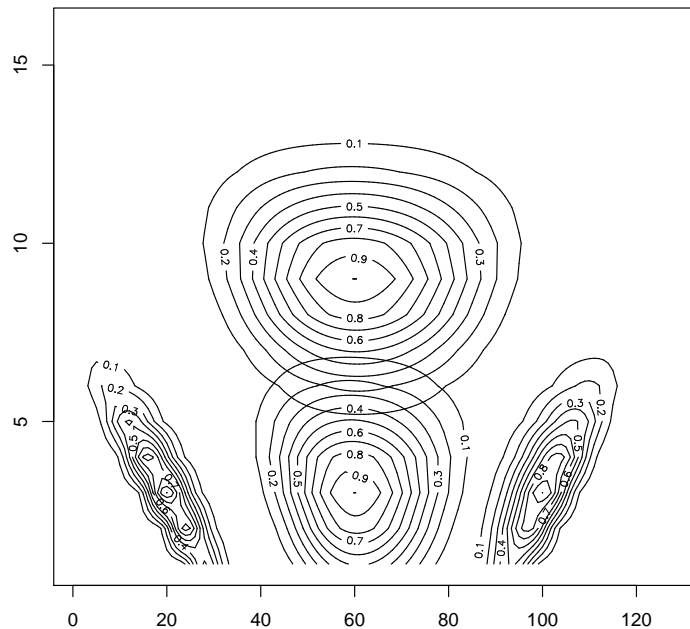
- A synthetic 1D example: 1D structure functions, original (dotted lines) and modelled with complex wavelets (bold).



- Variance tends to be underestimated in regions with large length scales. This has also been observed by other researchers.

# 1D Complex wavelets

- Because the covariances are complex (the real part at level 1 can be correlated to the imaginary part at level 2), the diagonal  $\mathbf{B}$  can still represent tilted structures, with some limitations.



# Non-periodic boundaries?

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- A single wavelet transform allowing symmetric boundaries, must be symmetric and odd length.
- (Anti-)symmetric wavelets can not be orthogonal (except the Haar wavelet).
- For the complex wavelets, symmetric wavelets are made possible by the combination of two inverse wavelets. The “errors” at the periodic boundaries thus compensate. BUT the first (smallest) scale needs a different wavelet.
- Kingsbury uses a symmetric, non-orthogonal wavelet. Others propose an orthogonal, even length, almost-symmetric filter, but this only allows periodic boundaries.



# Non-periodic: solution

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- Consider not the original domain  $X[1..N]$ , but the doubling  $X[1..N, N..1]$ , which is periodic by definition.
- On this domain, find an orthogonal almost-symmetric & odd length wavelet for the first stage. Such a wavelet existed in literature, but was hard to find.
- Use Kingsbury's wavelet on the next stages.
- Elements  $1..N$  are the real part, the second half are the corresponding imaginary parts.
- So the resulting solution is no longer a combination of 2 orthogonal wavelets, but 1 transform on the doubled domain!

# 2D complex wavelets

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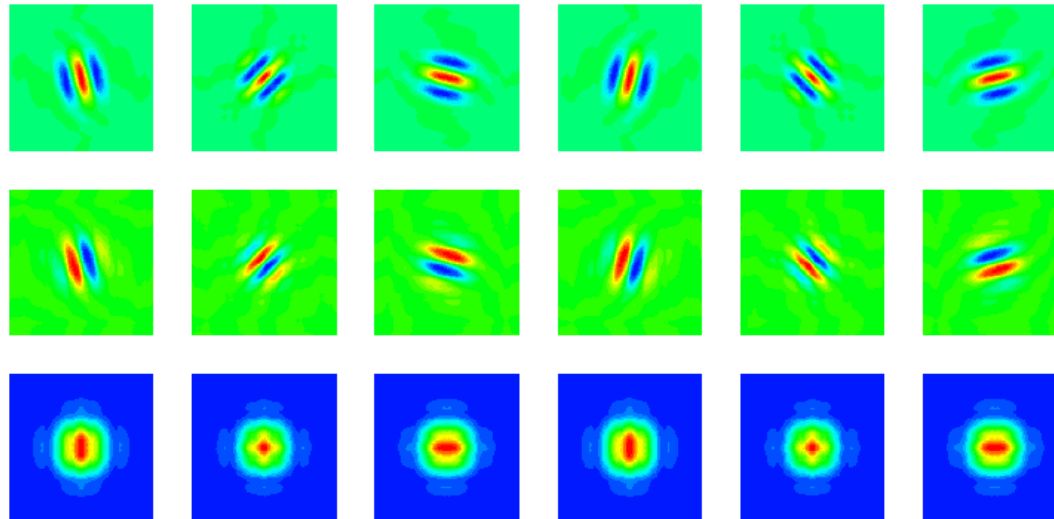
- Tensor products of the 2 dual-tree wavelets yields 4 different orthonormal 2D wavelet transforms, each of which has 3 different sectors.
- By taking certain linear combinations of these wavelet functions, we get a new (non-orthogonal) set of complex wavelets:

$$\begin{aligned}\Psi_l^1 &= (\psi_l^{11} + \psi_l^{22}) + i(\psi_l^{12} - \psi_l^{21}), \\ \Psi_l^2 &= (\psi_l^{12} + \psi_l^{21}) + i(\psi_l^{11} - \psi_l^{22}).\end{aligned}$$

# 2D complex wavelets – orientations

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- Real part, imaginary part and modulus of the 2D wavelets:

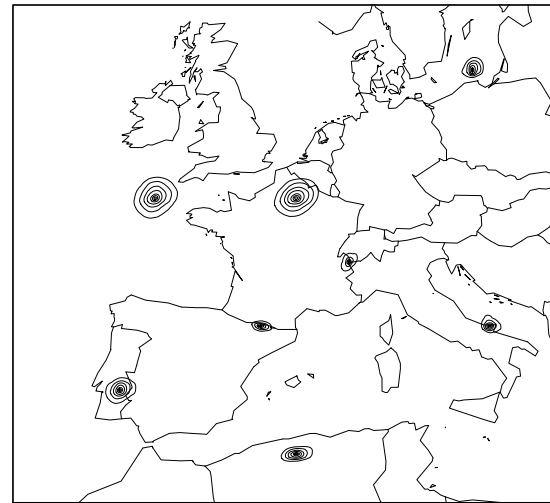
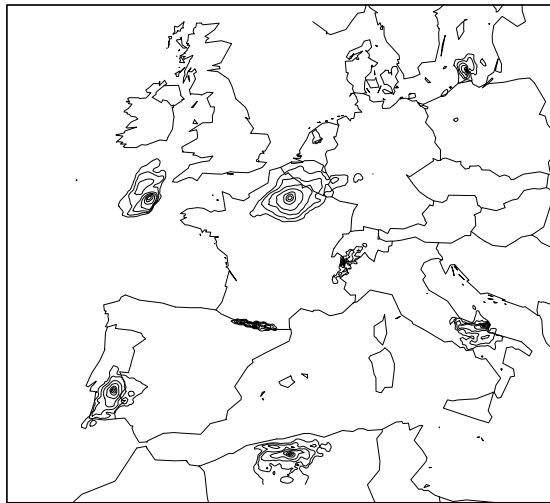


- → 6 distinct directional components at every scale!

# Complex wavelet transform

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- If we use these directional wavelets to diagonalise  $B$ :



# 3-dimensional covariances

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- We now move on to an  $N \times N \times L$  grid.
- **Vertical** covariances are best treated in a different manner: e.g. they are not evenly spaced. Also,  $L$  is usually much smaller than  $N$  (e.g. 46).
- In the standard approach,  $\mathbf{B}$  becomes **block-diagonal**: for every 2D component (spectral or wavelet) we get a “vertical”  $L \times L$  block.
- In the assimilation code, the eigenvectors of these blocks are computed numerically.

# Reducing memory cost

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- For 1 field, we still have  $4N^2 \times L^2$  components, which is too much in an operational setting.
- In ALADIN, the spectral components are also averaged by (total) wave number. So in 2D, there are only  $O(N)$  components left.
- We introduced 2 reductions.
- In the horizontal components, we may eliminate the smallest scales.

# Reducing the vertical matrices

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- For every scale and orientation (denoted  $l$ ) we calculate the *mean vertical matrix*  $\overline{\mathbf{B}}_l$ .
- This matrix can be rewritten in its eigenbasis  $\overline{\mathbf{E}}_l$ :

$$\overline{\mathbf{B}}_l = \overline{\mathbf{E}}_l \overline{\Lambda}_l \overline{\mathbf{E}}_l^*$$

- We then assume that the basis of eigenvectors  $\overline{\mathbf{E}}_l$  is representative for all local eigenbases. Thus we write for location  $i$ :

$$\mathbf{B}_{l,i} \approx \overline{\mathbf{E}}_l \Lambda_{l,i} \overline{\mathbf{E}}_l^*,$$

where  $\Lambda_{l,i} = \text{diag}(\overline{\mathbf{E}}_l^* \mathbf{B}_{l,i} \overline{\mathbf{E}}_l)$ .

# Reducing the vertical matrices

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- This averaging of vertical components has a rather strong smoothing effect. Some detail (especially close to the surface) is lost.
- This can be (partially) solved by taking out the variance before the reduction, thus only reducing the **correlations**, which are usually smoother.



# Implementation in ALADIN

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- Full wavelet treatment of **unbalanced part** of  $B$ . The statistical balance is still in Fourier space.
- Currently, only periodic wavelet version coded.
- Can be turned on with the switch LJBWAVELET.
- Still **very experimental**.
- $B_{wav}$  is calculated offline using a wavelet version of FESTAT.

# Implementation in ALADIN

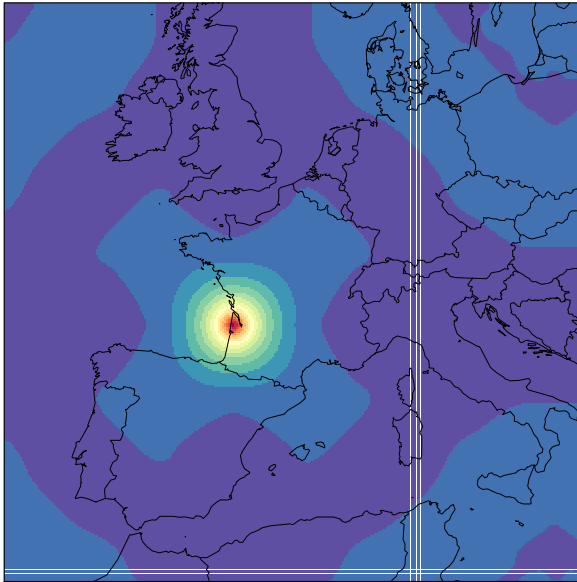
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- Wavelet domain must contain enough powers of 2 (so extend 300  $\rightarrow$  320)
- For the balance part, we go back to Fourier space.
- This requires bi-periodicisation at every iteration.
- Not only expensive, also no simple adjoint.
- Still causes noise unless we use the same (larger) extension zone for Fourier space as well.

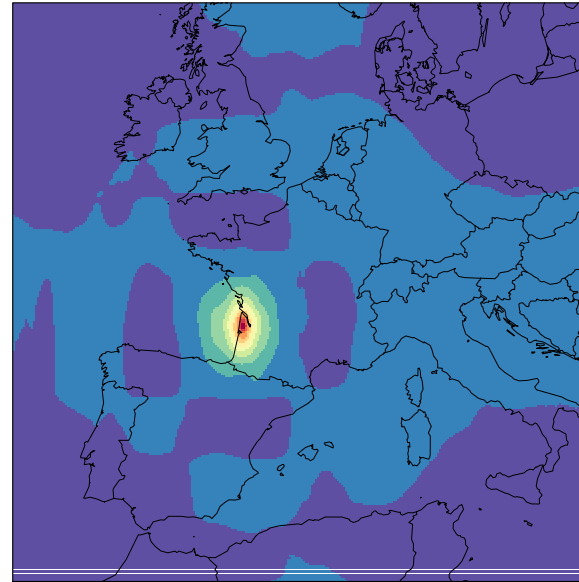
# Single T obs at 1000hPa

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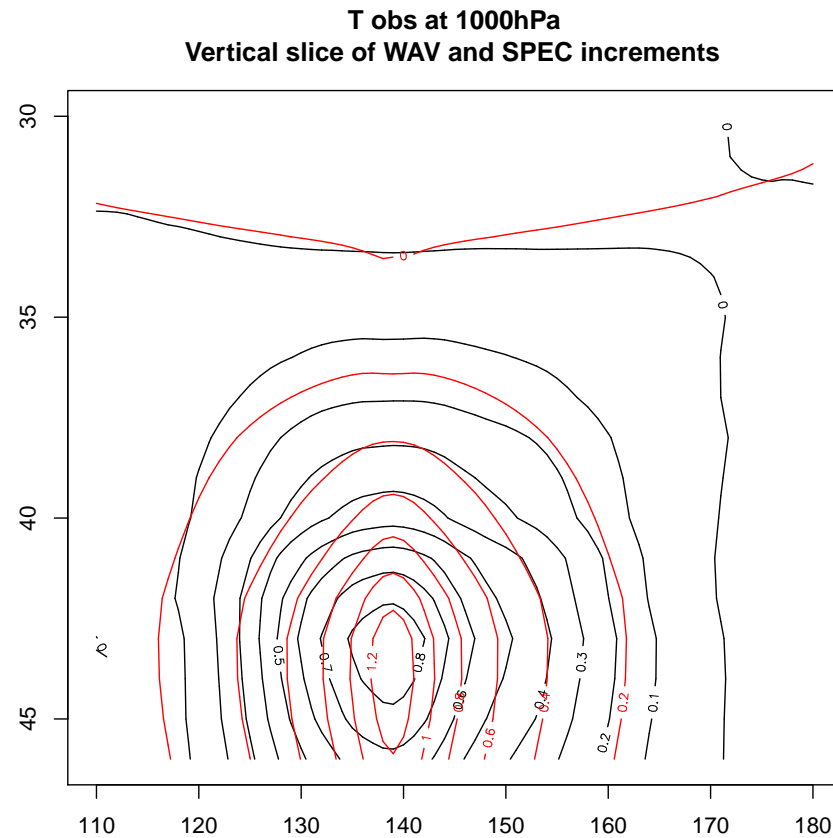
T obs at 1000 hPa  
T increment at lev 46  
B-SPEC



T obs at 1000 hPa  
T increment at lev 46  
B-WAV



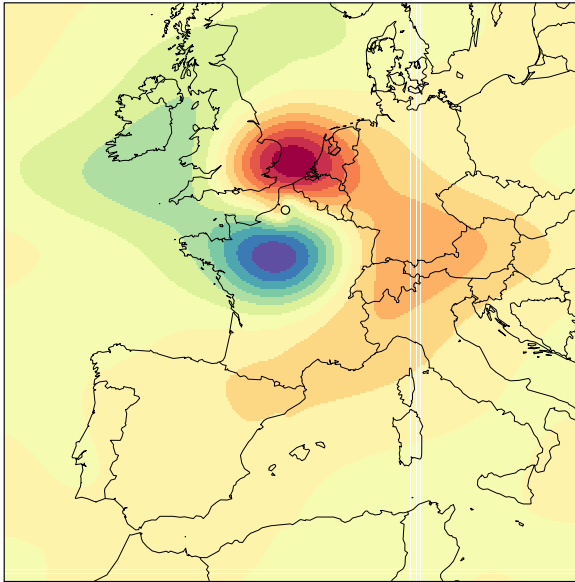
# Single T obs at 1000hPa



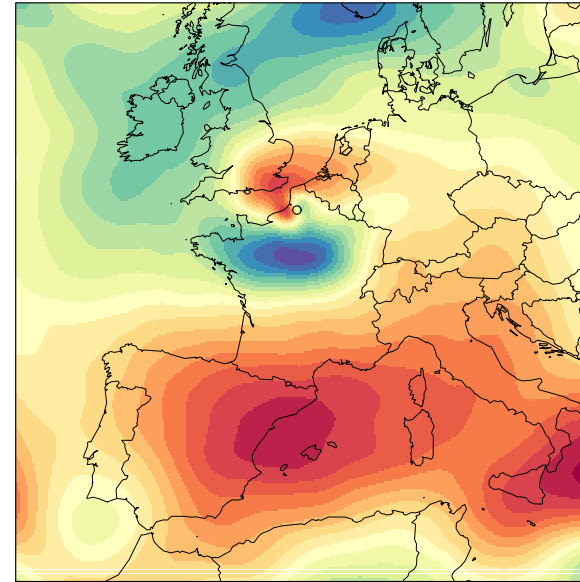
# Problem with wind increment?

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T obs at 300 hPa  
U increment lev 20  
B-SPEC



T obs at 300 hPa  
U increment lev 20  
B-WAV



# Conclusions: advantages

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- Using wavelet transform we can simplify the  $B$  matrix while maintaining its basic local and spatial (scale) features.
- Compared to standard ODWT, dual-tree complex wavelets have several advantages:
  - Approximate shift invariance (and hence reduced “artifacts”).
  - Improved directional resolution,
  - possibility of symmetric boundaries,
  - A phase that allows for tilted structure functions in 3D... But maybe it is removed by the vertical approximations...

# To do

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- Find out what is happening with the wind increments.
- Implement non-periodic transforms in ALADIN.
- Much more experimenting.
- Solve remaining noise issues & normalisation of variances.
- Better vertical representation (tilt,...)?